

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110421\
 Data File : BM032883.D
 Acq On : 04 Nov 2021 20:00
 Operator : CG/JU
 Sample : M4497-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RO-COMP-1

Quant Time: Nov 07 08:52:13 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 02 16:07:18 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.484	152	108152	20.000	ng	0.00
21) Naphthalene-d8	10.266	136	415553	20.000	ng	# 0.00
39) Acenaphthene-d10	14.148	164	236831	20.000	ng	0.00
64) Phenanthrene-d10	16.883	188	445137	20.000	ng	-0.01
76) Chrysene-d12	21.065	240	413225	20.000	ng	-0.01
86) Perylene-d12	23.177	264	436640	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.084	112	908368	146.427	ng	0.00
7) Phenol-d6	6.695	99	1024141	110.858	ng	0.00
23) Nitrobenzene-d5	8.642	82	756017	99.576	ng	0.00
42) 2,4,6-Tribromophenol	15.642	330	346432	130.613	ng	0.00
45) 2-Fluorobiphenyl	12.760	172	1573770	103.065	ng	0.00
79) Terphenyl-d14	19.512	244	1910203	100.925	ng	-0.01
Target Compounds						
32) Benzoic acid	9.595	122	16604	3.674	ng	# 81

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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